CLAIMS:

1. The use of a compound of Formula (I):

Formula (I)

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wherein:

X and **Y** are independently selected from: oxygen, sulphur and $(-CR^aR^b-)_n$; wherein: **n** is an integer of from 1 to 4 and

 $\mathbf{R}^{\mathbf{a}}$ and $\mathbf{R}^{\mathbf{b}}$ are each independently selected from hydrogen, C_{1-6} alkyl, C_{1-6} alkoxy, halo, hydroxy, C_{1-6} alkanoyloxy, C_{3-12} cycloalkyl and optionally substituted phenyl or $\mathbf{R}^{\mathbf{a}}$ and $\mathbf{R}^{\mathbf{b}}$ together form a C_{5-12} spirocycloalkyl or a carbonyl; with the proviso that at least one of \mathbf{X} and \mathbf{Y} is $(-C\mathbf{R}^{\mathbf{a}}\mathbf{R}^{\mathbf{b}}-)_{\mathbf{n}}$ and with the further proviso that when \mathbf{X} and \mathbf{Y} are both $(-C\mathbf{R}^{\mathbf{a}}\mathbf{R}^{\mathbf{b}}-)_{\mathbf{n}}$ and $\mathbf{R}^{\mathbf{a}}$ are hydrogen and \mathbf{n} is 1, then $\mathbf{R}^{\mathbf{1}}$ and $\mathbf{R}^{\mathbf{3}}$ are both aryl;

15 \mathbb{R}^2 is hydrogen, a C_{1-8} alkyl or benzyl;

 ${\bf R}^1$ and ${\bf R}^3$ are independently selected from

- (a) phenyl or phenoxy wherein the phenyl or phenoxy group is optionally substituted with 1 to 5 substituents independently selected from phenyl, C₁₋₆alkyl, C₁₋₆alkoxy, phenoxy, hydroxy, fluorine, chlorine, bromine, nitro, trifluoromethyl, carboxy, C₁₋₄alkoxycarbony and -(CH₂)_pNR₄R₅ wherein p is 0 or 1, and R⁴ and R⁵ are independently selected from hydrogen and C₁₋₄alkyl;
- (b) naphth-1-yl or naphth-2-yl wherein the naphthyl group is optionall substituted with from 1 to 3 substituents independently selected from phenyl, C₁₋₆alkyl, C₁₋₆alkoxy, phenoxy, hydroxy, fluorine, chlorine, bromine, nitro, trifluoromethyl, carboxy, C₁₋₄alkoxycarbony and -(CH₂)_pNR₄R₅ wherein p, R⁴ and R⁵ are as defined above;
- (c) arylC₁₋₆alkyl;
- (d) C₁₋₂₀alkyl or C₁₋₂₀alkenyl; and
- (e) adamantyl or a C₃₋₁₂cycloalkyl;
- or a pharmaceutically acceptable salt, pro-drug or solvate thereof in the manufacture of a medicament for the treatment of diabetes and/or obesity.

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- 2. The use according to claim 1 wherein \mathbb{R}^1 is phenyl.
- 3. The use according to claim 2 wherein \mathbb{R}^1 is phenyl disubstituted in the 2,6-positions.
- 4. The use according to any one of claims 1, 2 or 3 wherein \mathbb{R}^3 is phenyl.
- 5. The use according to claim 4 wherein \mathbb{R}^3 is phenyl disubstituted in the 2,6-positions.
- 10 6. The use according to claim 1 wherein \mathbb{R}^1 is phenyl disubstituted in the 2,6-positions and \mathbb{R}^3 is phenyl trisubstituted in the 2,4,6-positions.
 - 7. The use according to claim 1 wherein \mathbb{R}^1 is 2,6-bis(1-methylethyl)phenyl and \mathbb{R}^3 is 2,6-bis(1-methylethyl)phenyl or 2,4,6-tris(1-methylethyl)phenyl.
 - 8. The use according to claim 1 wherein: one of $\mathbf{R^1}$ and $\mathbf{R^3}$ is the group

$$R_1^6$$
 — $(CH_2)_t$ — C — $(CH_2)_w$ — R^8

wherein

20 **t** is 0 to 4;

w is 0 to 4 with the proviso that the sum of t and w is not greater than 5;

- ${f R}^6$ and ${f R}^7$ are independently selected from hydrogen or $C_{1\text{-}6}$ alkyl, or when ${f R}^6$ is hydrogen, ${f R}^7$ can be selected from the groups defined for ${f R}^8$; and ${f R}^8$ is phenyl optionally substituted with from 1 to 3 substituents selected $C_{1\text{-}6}$ alkyl $C_{1\text{-}6}$ alkoxy, phenoxy, hydroxy, fluorine, chlorine, bromine, nitro, trifluoromethyl, carboxy, $C_{1\text{-}4}$ alkoxycarbonyl, or -(CH₂)pNR⁴R⁵ wherein p, R⁴ and R⁵ are as defined above.
- 9. The use according to claim 1 wherein
- 30 X is oxygen;

Y is (CR^aR^b)_n wherein

n is an integer of from 1 to 4 and

 $\mathbf{R}^{\mathbf{a}}$ and $\mathbf{R}^{\mathbf{b}}$ are each independently hydrogen, C_{1-6} alkyl, optionally substituted phenyl, halo, hydroxy, C_{1-6} alkoxy, C_{1-6} alkanoyloxy, cycloalkyl, or $\mathbf{R}^{\mathbf{a}}$ and $\mathbf{R}^{\mathbf{b}}$ taken together form a carbonyl or C_{3-10} spirocycloalkyl;

 \mathbb{R}^1 is selected from optionally substituted phenyl, C_{1-10} alkyl or C_{3-10} cycloalkyl;

R² is hydrogen;

 ${\bf R^3}$ is selected from optionally substituted phenyl, C_{1-10} alkyl, C_{3-8} cycloalkyl and optionally substituted phenoxy.

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10. The use according to claim 1 wherein

X is oxygen;

Y is $(CR^aR^b)_n$ wherein n is an integer of from 1 to 2;

R¹ is optionally substituted phenyl;

15 R² is hydrogen;

 ${f R}^3$ is selected from optionally substituted phenyl, optionally substituted phenoxy, C_{1-10} alkyl, and C_{3-10} cycloalkyl;

 ${f R}^a$ and ${f R}^b$ are independently selected from hydrogen, $C_{1\text{-}6}$ alkyl, optionally substituted phenyl, halogen, hydroxy, $C_{1\text{-}6}$ alkoxy, $C_{1\text{-}6}$ alkanoyloxy, cycloalkyl, or ${f R}^a$ and ${f R}^b$ taken together form a carbonyl or a spirocycloalkyl.

11. The use according to claim 1 wherein

X is oxygen;

Y is (-CR^aR^b-)n wherein n is an integer of from 1 to 4 and R' and R" are each independently hydrogen, alkyl, alkoxy, halogen, hydroxy, acyloxy, cycloalkyl, phenyl optionally substituted or R' and R" together form a spirocycloalkyl or a carbonyl;

 $\boldsymbol{R^1}$ and $\boldsymbol{R^3}$ are independently selected from

- (a) phenyl or phenoxy wherein the phenyl or phenoxy group is optionally substituted with 1 to 5 substituents independently selected from
- phenyl, C_{1-6} alkyl, C_{1-6} alkoxy, phenoxy, hydroxy, fluorine, chlorine, bromine, nitro, trifluoromethyl, carboxy, C_{1-4} alkoxycarbony and $-(CH_2)_pNR_4R_5$ wherein p is 0 or 1, and R^4 and R^5 are independently selected from hydrogen or C_{1-4} alkyl;

- (b) naphth-1-yl or naphth-2-yl wherein the naphthyl group is optionally substituted with from 1 to 3 substituents independently selected from
- phenyl, C_{1-6} alkyl, C_{1-6} alkoxy, phenoxy, hydroxy, fluorine, chlorine, bromine, nitro, trifluoromethyl, carboxy, C_{1-4} alkoxycarbony and- $(CH_2)_pNR_4R_5$ wherein p, R^4 and R^5 are as defined above;
- (c) arylC₁₋₆alkyl;

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- (d) C₁₋₂₀alkyl or C₁₋₂₀alkenyl; and
- (e) adamantyl or a C₃₋₁₂cycloalkyl

 \mathbb{R}^2 is hydrogen, a \mathbb{C}_{1-8} alkyl or benzyl;

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- 12. The use according to claim 1 wherein the compound is selected from:
 - Sulfamic acid (phenylacetyl)-2,6-bis(1-methylethyl)phenyl ester,
 - Sulfamic acid[[2,6-bis(1-methylethyl)phenyl]acetyl]-2,6-bis(1-methylethyl)phenyl ester,
- Sulfamic acid [[2,4,6-tris(1-methylethyl)phenyl]acetyl-2,4,6-tris(1-methylethyl)phenyl ester,
 - Sulfamic acid[[2,6-bis(1-methylethyl)phenyl]acetyl]-2,4,6-tris(1-methylethyl)phenyl ester,
 - Sulfamic acid[adamantaneacetyl]-2,6-bis[1-methylethyl)phenyl ester,
- Sulfamic acid[[2,6-bis(1-methylethyl)phenyl]acetyl]-2,6-bis(1-methylethyl)phenyl ester-sodium salt,
 - Sulfamic acid[[2,4,6-tris(1-methylethyl)phenyl]acetyl]-2,6-bis(1-methylethyl)phenyl ester-sodium salt,
 - Sulfamic acid (decanoyl)-2,6-bis-(1-methylethyl)phenyl ester,
- Sulfamic acid (dodecanoyl)-2,6-bis-(1-methylethyl)phenyl ester,
 - 2. 6-Bis(1-methylethyl)-N-[[[2,4,6-tris(1-methylethyl)phenyl]methyl]sulfonyl] benzeneacetamide,
 - 2,6-Bis(1-methylethyl)-N-[[[2,4,6-tris(1-methylethyl)phenyl]methyl]sulfonyl] benzeneacetamide-sodium salt.

- 2,6-Bis(1-methylethyl)phenyl[[[2,4,6-tris(1-methylethyl)phenyl]methyl]sulfonyl]carbamate,
- 2,6-Bis(1-methylethyl)phenyl[[[2,4,6-tris(1-methylethyl)phenyl]methyl]sulfonyl]carbamate-sodium salt,
- Sulfamic acid (1-oxo-3,3-diphenylpropyl)-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid [2,6-dichlorophenyl(acetyl)]-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid [2,6-dichlorophenyl(acetyl)]-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid trans-[(2-phenylcyclopropyl)carbonyl]-2,6-bis(1-methylethyl)phenyl ester,
- Sulfamic acid [2,5-dimethoxyphenyl(acetyl)]-2,6-bis(1-methylethyl)phenyl ester,
 Sulfamic acid [2,4,6-trimethoxyphenyl(acetyl)]-2,6-bis(1-methylethyl)phenyl ester,
 Sulfamic acid [2,4,6-trimethylphenyl(acetyl)]-2,6-bis(1-methylethyl)phenyl ester,
 Sulfamic acid [2-thiophenyl(acetyl)]-2,6-bis(1-methylethyl)phenyl ester,
 Sulfamic acid [3-thiophenyl(acetyl)]-2,6-bis(1-methylethyl)phenyl ester,
- Sulfamic acid [2-methoxyphenyl(acetyl)]-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid (oxophenylacetyl)-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid [2-trifluoromethylphenyl(acetyl)]-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid (1-oxo-2-phenylpropyl)-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid (cyclopentylphenylacetyl)-2,6-bis(1-methylethyl)phenyl ester,
- Sulfamic acid (cyclopentylphenylacetyl)-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid (cyclohexylacetyl)-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid (diphenylacetyl)-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid (triphenylacetyl)-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid [(11-phenylcyclopentyl)carbonyl]-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid (3-methyl-1-oxo-2-phenylpentyl)-2,6-bis(1-methylethyl)phenyl ester,
- Sulfamic acid (1-oxo-2-phenylbutyl)-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid (cyclohexylphenylacetyl)-2,6-bis(1-methylethyl)phenyl ester,

 Sulfamic acid (1-oxo-2,2-diphenylpropyl)-2,6-bis(1-methylethyl)phenyl ester,

- Sulfamic acid [(9H-fluoren-9-yl)carbonyl]-2,6-bis(1-methylethyl)phenyl ester,
- Sulfamic acid (1-oxo-3-phenylpropyl)-2,6-bis(1-methylethyl)phenyl ester,
- Sulfamic acid [1-oxo-3-[2,4,6-tris(1-methylethyl)phenyl-2-propenyl]-2,6-bis(1-methylethy l)phenyl ester,
- Sulfamic acid [1-oxo-3-[2,4,6-tris(1-methylethyl)phenyl]propyl]-2,6-bis(1-methylethyl)ph enyl ester,
 - Sulfamic acid [(acetyloxy)[2,4,6-tris(1-methylethyl)phenyl]acetyl]-2,6-bis(1-methylethyl)phenyl ester,
 - Sulfamic acid [hydroxy[2,4,6-tris(1-methylethyl)phenyl]acetyl]-2,6-bis(1-methylethyl)phe nyl ester,
 - Sulfamic acid [fluoro[2,4,6-tris(1-methylethyl)phenyl]acetyl]-2,6-bis(1-methylethyl)phen yl ester,
 - Sulfamic acid (3-methyl-1-oxo-2-phenylpentyl)-2,6-bis(1-methylethyl)phenyl ester sodium salt,
- Sulfamic acid [[2,4,6-tris(1-methylethyl)phenoxy]acetyl]-2,6-bis(1-methylethyl)phenyl ester,
 - Sulfamic acid [[2,6-bis(1-methylethyl)phenoxy]acetyl]-2,6-bis(1-methylethyl)phenyl ester, and
 - Sulfamic acid [[2,4,6-tris(1-methylethyl)phenyl]acetyl]-2,6-bis(phenyl)phenyl ester.
- or pharmaceutically acceptable salt, pro-drug or solvate.

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- 13. The use according to claim 1 wherein the compound is: sulfamic acid[[2,4,6-tris(1-methylethyl)phenyl]acetyl-2,6-bis(1-methylethyl)phenyl ester.
- or pharmaceutically acceptable salt, pro-drug or solvate.
 - 14. The use according to any one of the preceding claims wherein the use is the manufacture of a medicament for the treatment of type II diabetes.

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- 15. The use according to any one of the preceding claims wherein the use is the manufacture of a medicament for the treatment of obesity.
- 16. The use according to any one of the preceding claims wherein the use is themanufacture of a medicament for the treatment of insulin resistance.
 - 17. The use according to any one of the preceding claims wherein the use is the manufacture of a medicament for the treatment of impaired glucose tolerance.